Introduction to Machine Learning

Gaussian Processes

$f(x)$

$$\sim \mathcal{N}\left(\mu, \Sigma\right)$$

Learning goals

- GPs model distributions over functions
- The marginalization property makes this distribution easily tractable
- GPs are fully specified by mean and covariance function
- GPs are indexed families

WEIGHT-SPACE VIEW

- Until now we considered a hypothesis space $\mathcal{H}$ of parameterized functions $f(\mathbf{x} \mid \boldsymbol{\theta})$ (in particular, the space of linear functions).
- Using Bayesian inference, we derived distributions for $\boldsymbol{\theta}$ after having observed data $\mathcal{D}$.
- Prior believes about the parameter are expressed via a prior distribution $q(\boldsymbol{\theta})$, which is updated according to Bayes' rule

$$\underbrace{p(\boldsymbol{\theta}|\mathbf{X}, \mathbf{y})}_{\text{posterior}} = \frac{\overbrace{p(\mathbf{y}|\mathbf{X}, \boldsymbol{\theta})}^{\text{likelihood}} \overbrace{q(\boldsymbol{\theta})}^{\text{prior}}}{\underbrace{p(\mathbf{y}|\mathbf{X})}_{\text{marginal}}}.$$
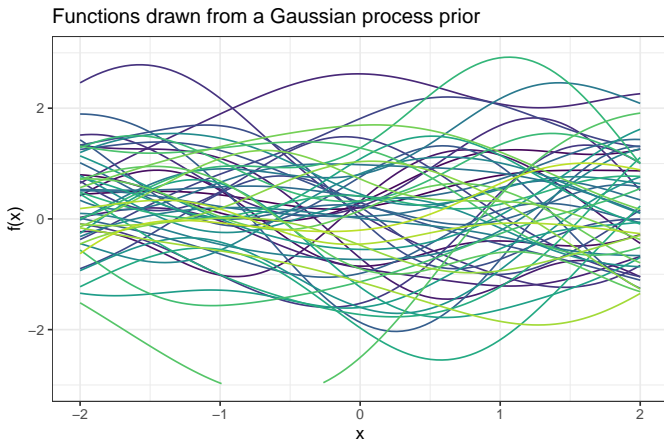
FUNCTION-SPACE VIEW

Let us change our point of view:

- Instead of "searching" for a parameter θ in the parameter space, we directly search in a space of "allowed" functions $\mathcal{H}$.
- We still use Bayesian inference, but instead specifying a prior distribution over a parameter, we specify a prior distribution **over functions** and update it according to the data points we have observed.
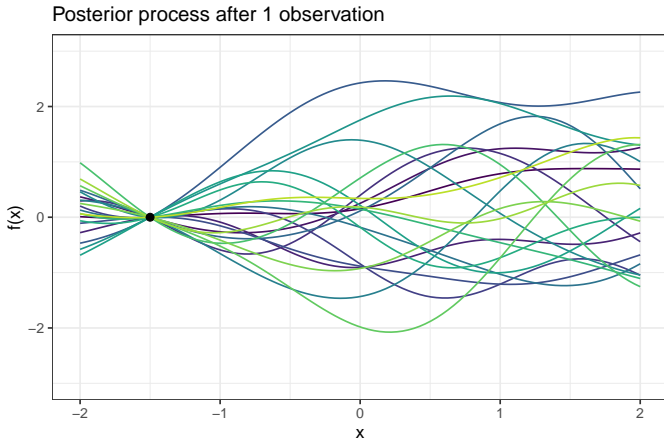
FUNCTION-SPACE VIEW

Intuitively, imagine we could draw a huge number of functions from some prior distribution over functions [*].

Functions drawn from a Gaussian process prior

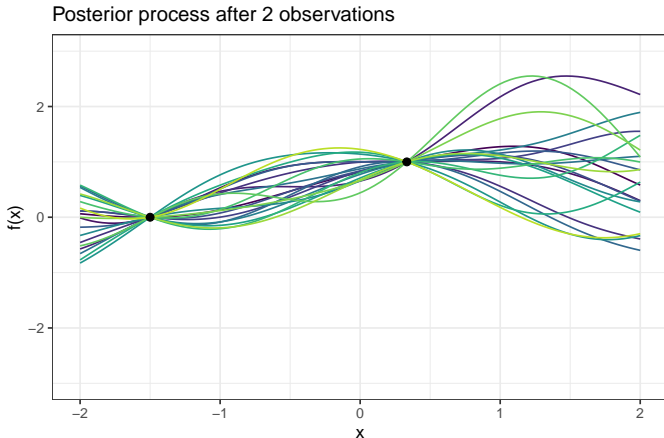[*] We will see in a minute how distributions over functions can be specified.

FUNCTION-SPACE VIEW

After observing some data points, we are only allowed to sample those
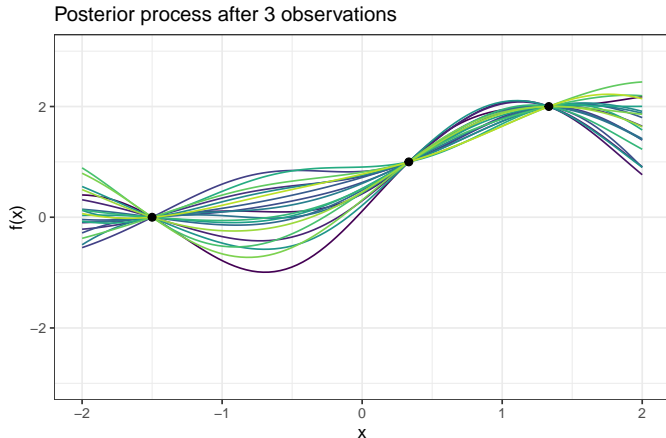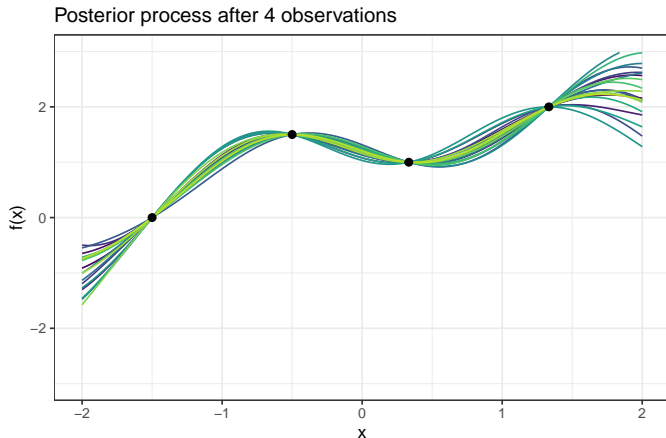functions, that are consistent with the data.

Posterior process after 1 observation

FUNCTION-SPACE VIEW

After observing some data points, we are only allowed to sample those functions, that are consistent with the data.

Posterior process after 2 observations

FUNCTION-SPACE VIEW

After observing some data points, we are only allowed to sample those functions, that are consistent with the data.

Posterior process after 3 observations

FUNCTION-SPACE VIEW

As we observe more and more data points, the variety of functions consistent with the data shrinks.

Posterior process after 4 observations

FUNCTION-SPACE VIEW

Inutitively, there is something like "mean" and a "variance" of a distribution over functions.

Posterior process after 4 observations
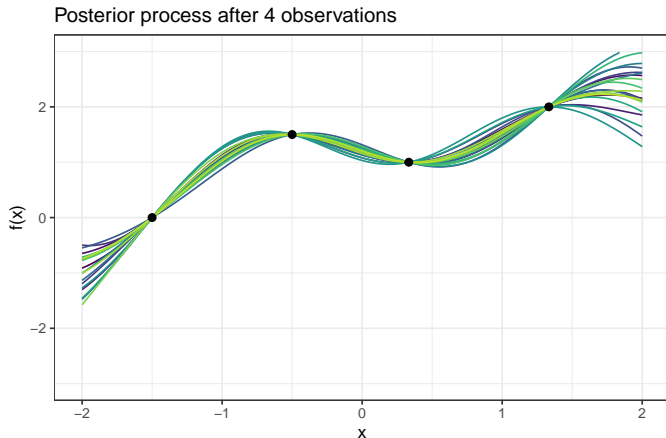
WEIGHT-SPACE VS. FUNCTION-SPACE VIEW

Weight-Space View	**Function-Space View**

Parameterize functions

Example: $f(\mathbf{x} \mid \boldsymbol{\theta}) = \boldsymbol{\theta}^\top \mathbf{x}$

Define distributions on $\boldsymbol{\theta}$	Define distributions on f

Inference in parameter space Θ	Inference in function space $\mathcal{H}$

Next, we will see how we can define distributions over functions mathematically.

Distributions on Functions

DISCRETE FUNCTIONS

For simplicity, let us consider functions with finite domains first.

Let $\mathcal{X} = \left\{ \mathbf{x}^{(1)}, \ldots, \mathbf{x}^{(n)} \right\}$ be a finite set of elements and $\mathcal{H}$ the set of all functions from $\mathcal{X} \rightarrow \mathbb{R}$.
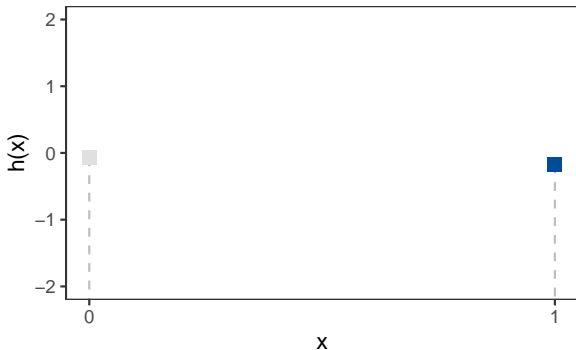
Since the domain of any $h(.) \in \mathcal{H}$ has only *n* elements, we can represent the function $h(.)$ compactly as a *n*-dimensional vector

$$\boldsymbol{h} = \left[h\left(\mathbf{x}^{(1)} \right), \ldots, h\left(\mathbf{x}^{(n)} \right) \right].$$

DISCRETE FUNCTIONS

Example 1: Let us consider $h : \mathcal{X} \to \mathcal{Y}$ where the input space consists of **two** points $\mathcal{X} = \{0, 1\}$.
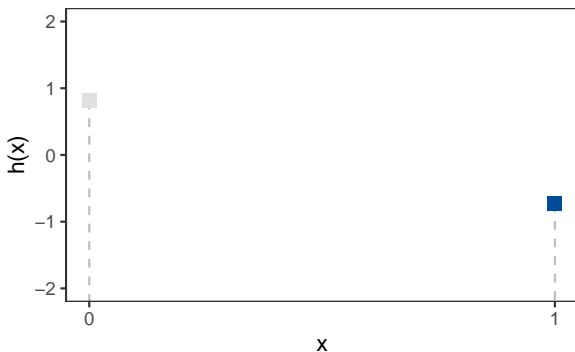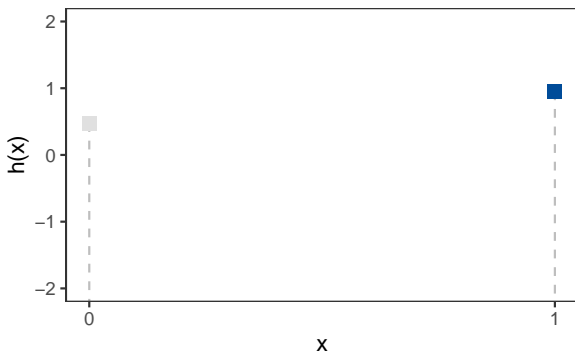
Examples for functions that live in this space:

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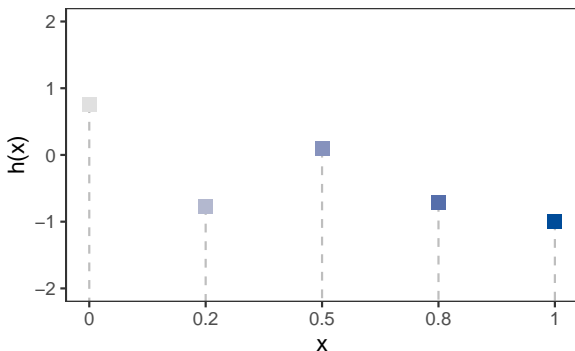
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Examples for functions that live in this space:

DISCRETE FUNCTIONS

Example 2: Let us consider $h : \mathcal{X} \to \mathcal{Y}$ where the input space consists of **five** points $\mathcal{X} = \{0, 0.25, 0.5, 0.75, 1\}$.
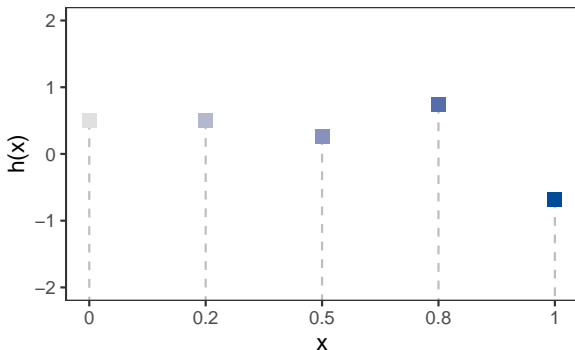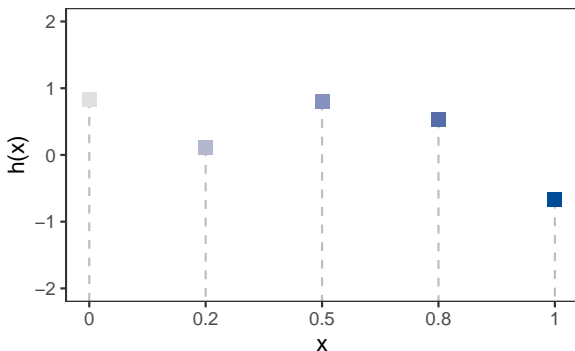
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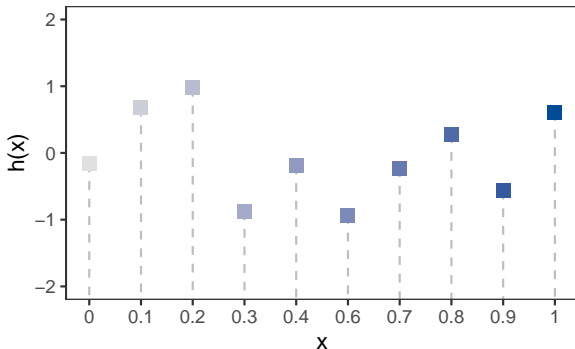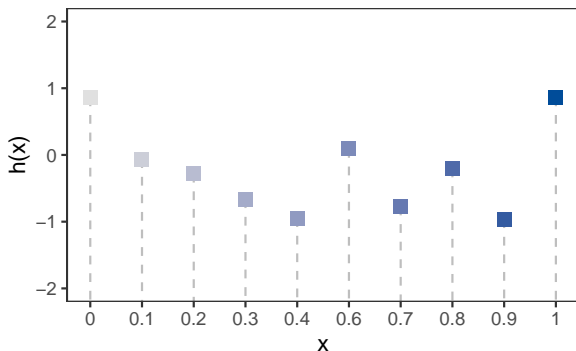
Examples for functions that live in this space:

DISCRETE FUNCTIONS

Example 3: Let us consider $h : \mathcal{X} \to \mathcal{Y}$ where the input space consists of **ten** points.

Examples for functions that live in this space:

DISCRETE FUNCTIONS

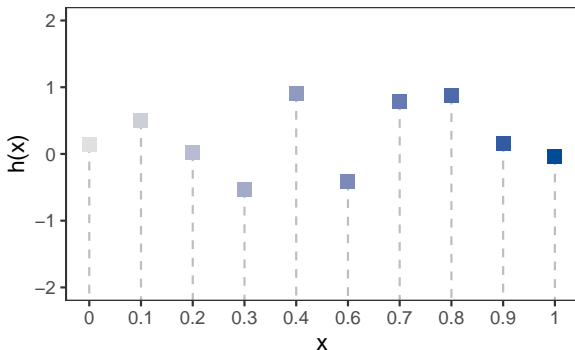**Example 3:** Let us consider $h : \mathcal{X} \to \mathcal{Y}$ where the input space consists of **ten** points.

Examples for functions that live in this space:

DISCRETE FUNCTIONS

Example 3: Let us consider $h : \mathcal{X} \to \mathcal{Y}$ where the input space consists of **ten** points.

Examples for functions that live in this space:

DISTRIBUTIONS ON DISCRETE FUNCTIONS

One natural way to specify a probability function on discrete function $h \in \mathcal{H}$ is to use the vector representation

$$\boldsymbol{h} = \left[h\left(\mathbf{x}^{(1)}\right), h\left(\mathbf{x}^{(2)}\right), \ldots, h\left(\mathbf{x}^{(n)}\right) \right]$$

of the function.

Let us see $\boldsymbol{h}$ as a n-dimensional random variable. We will further assume the following normal distribution:
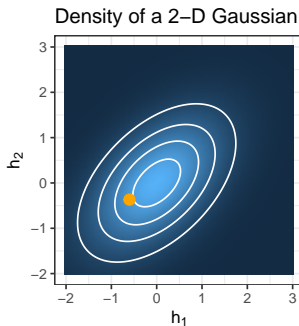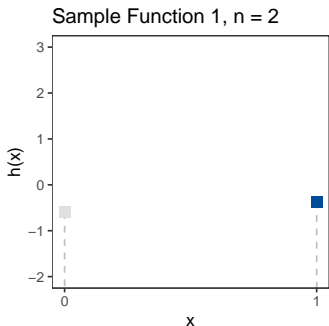
$$\boldsymbol{h} \sim \mathcal{N}\left(\boldsymbol{m}, \boldsymbol{K}\right).$$

Note: For now, we set $\boldsymbol{m} = \boldsymbol{0}$ and take the covariance matrix $\boldsymbol{K}$ as given. We will see later how they are chosen / estimated.

DISCRETE FUNCTIONS

Example 1 (continued): Let $h : \mathcal{X} \to \mathcal{Y}$ be a function that is defined on **two** points $\mathcal{X}$. We sample functions by sampling from a two-dimensional normal variable

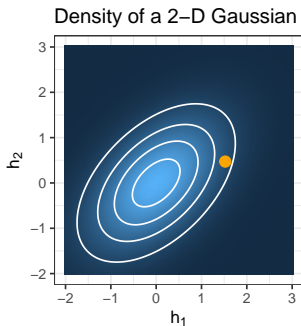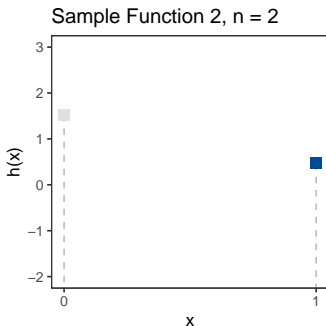$$\boldsymbol{h} = [h(1), h(2)] \sim \mathcal{N}(\boldsymbol{m}, \boldsymbol{K})$$

In this example, $m = (0, 0)$ and $K = \begin{pmatrix} 1 & 0.5 \\ 0.5 & 1 \end{pmatrix}$.

DISCRETE FUNCTIONS

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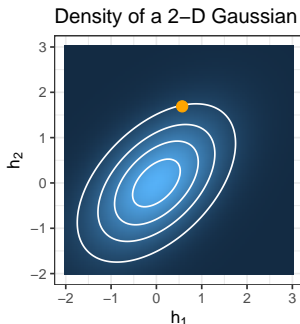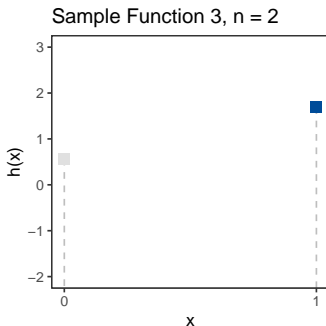$$\boldsymbol{h} = [h(1), h(2)] \sim \mathcal{N}(\boldsymbol{m}, \boldsymbol{K})$$

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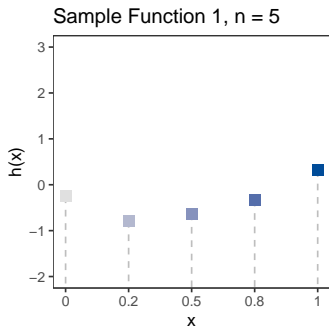$$\boldsymbol{h} = [h(1), h(2)] \sim \mathcal{N}(\boldsymbol{m}, \boldsymbol{K})$$

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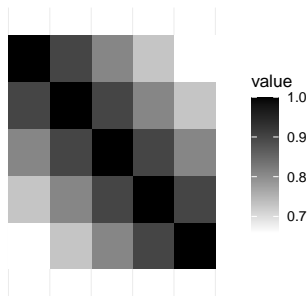$$\boldsymbol{h} = [h(1), h(2), h(3), h(4), h(5)] \sim \mathcal{N}(\boldsymbol{m}, \boldsymbol{K})$$

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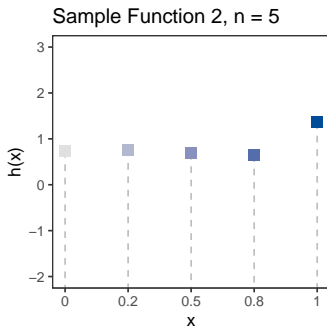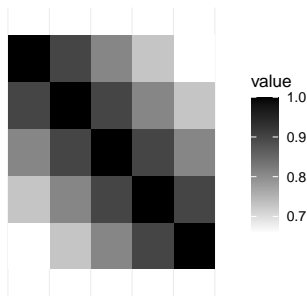## DISCRETE FUNCTIONS

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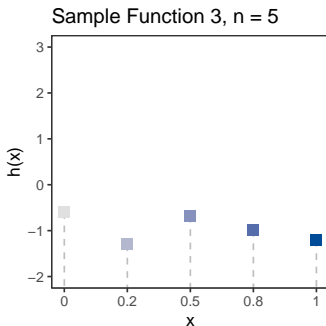
DISCRETE FUNCTIONS

Example 3 (continued): Let us consider $h : \mathcal{X} \to \mathcal{Y}$ where the input space consists of **ten** points. We sample functions by sampling from ten-dimensional normal variable

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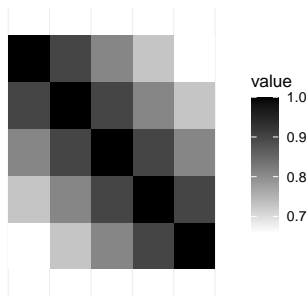# DISCRETE FUNCTIONS

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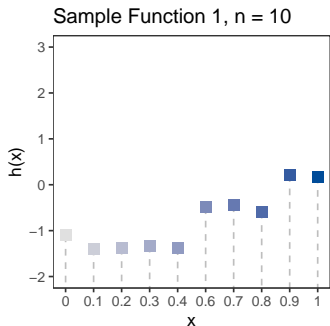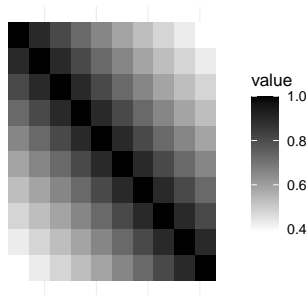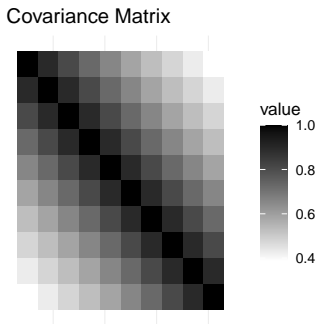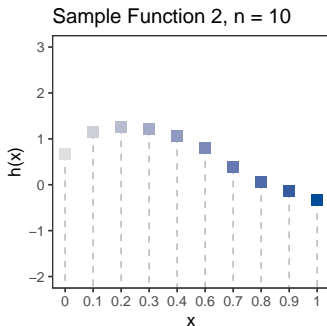## DISCRETE FUNCTIONS

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$$\boldsymbol{h} = [h(1), h(2), \ldots, h(10)] \sim \mathcal{N}(\boldsymbol{m}, \boldsymbol{K})$$

ROLE OF THE COVARIANCE FUNCTION

Note that the covariance controls the "shape" of the drawn function.
Consider two extreme cases where function values are

a) strongly correlated: $K = \begin{pmatrix} 1 & 0.99 & \ldots & 0.99 \\ 0.99 & 1 & \ldots & 0.99 \\ & & \ddots & \\ 0.99 & 0.99 & & 0.99 \\ 0.99 & \ldots & 0.99 & 1 \end{pmatrix}$

b) uncorrelated: $K = I$
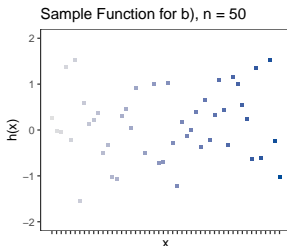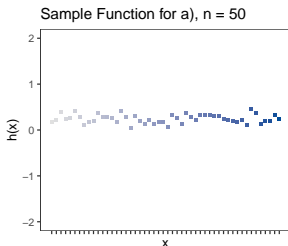
ROLE OF THE COVARIANCE FUNCTION

- "Meaningful" functions (on a numeric space $\mathcal{X}$) may be characterized by a spatial property:

    If two points $\mathbf{x}^{(i)}, \mathbf{x}^{(j)}$ are close in $\mathcal{X}$-space, their function values $f(\mathbf{x}^{(i)}), f(\mathbf{x}^{(j)})$ should be close in $\mathcal{Y}$-space.

    In other words: If they are close in $\mathcal{X}$-space, their functions values should be **correlated**!

- We can enforce that by choosing a covariance function with

    $$\mathbf{K}_{ij} \text{ high, if } \mathbf{x}^{(i)}, \mathbf{x}^{(j)} \text{ close.}$$

ROLE OF THE COVARIANCE FUNCTION

- We can compute the entries of the covariance matrix by a function that is based on the distance between $\mathbf{x}^{(i)}, \mathbf{x}^{(j)}$, for example:

  c) Spatial correlation: $K_{ij} = k(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}) = \exp\left(-\frac{1}{2}\left|\mathbf{x}^{(i)} - \mathbf{x}^{(j)}\right|^2\right)$

Sample Function for b) K = I, n = 50   Sample Function for c), n = 50
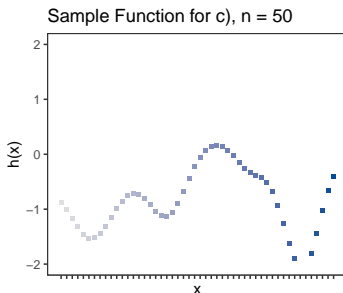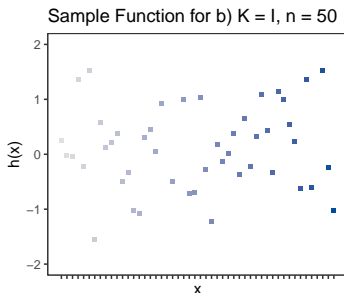
Note: $k(\cdot, \cdot)$ is known as the **covariance function** or **kernel**. It will be studied in more detail later on.

Gaussian Processes

FROM DISCRETE TO CONTINUOUS FUNCTIONS

- We defined distributions on functions with discrete domain by defining a Gaussian on the vector of the respective function values

$$\mathbf{h} = [h(\mathbf{x}^{(1)}), h(\mathbf{x}^{(2)}), \ldots, h(\mathbf{x}^{(n)})] \sim \mathcal{N}(\boldsymbol{m}, \boldsymbol{K})$$

- We can do this for $n \to \infty$ (as "granular" as we want)
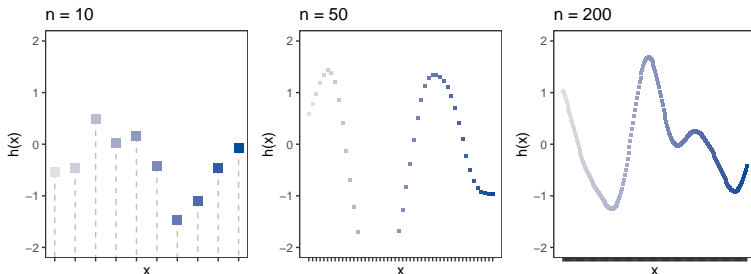
FROM DISCRETE TO CONTINUOUS FUNCTIONS

- No matter how large *n* is, we are still considering a function over a discrete domain.
- How can we extend our definition to functions with **continuous domain** $\mathcal{X} \subset \mathbb{R}$?

GAUSSIAN PROCESSES: INTUITION

- Intuitively, a function *f* drawn from **Gaussian process** can be understood as an "infinite" long Gaussian random vector.
- It is unclear how to handle an "infinite" long Gaussian random vector!

?

GAUSSIAN PROCESSES: INTUITION

- Thus, it is required that for **any finite set** of inputs
  $\{\mathbf{x}^{(1)}, \ldots, \mathbf{x}^{(n)}\} \subset \mathcal{X}$, the vector **f** has a Gaussian distribution

  $$\mathbf{f} = \left[f\left(\mathbf{x}^{(1)}\right), \ldots, f\left(\mathbf{x}^{(n)}\right) \right] \sim \mathcal{N}\left(\mathbf{m}, \mathbf{K}\right),$$

  with $\mathbf{m}$ and $\mathbf{K}$ being calculated by a mean function $m(.)$ /
  covariance function $k(.,.)$.

- This property is called **Marginalization Property**.

Sample Function, n = 5

$f(x)$

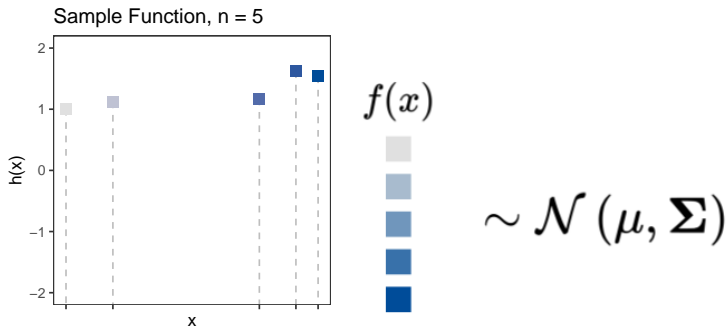$\sim \mathcal{N}\left(\mu, \mathbf{\Sigma}\right)$

GAUSSIAN PROCESSES: INTUITION

- Thus, it is required that for **any finite set** of inputs $\{\mathbf{x}^{(1)}, \ldots, \mathbf{x}^{(n)}\} \subset \mathcal{X}$, the vector **f** has a Gaussian distribution

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Sample Function, n = 10

$f(x)$

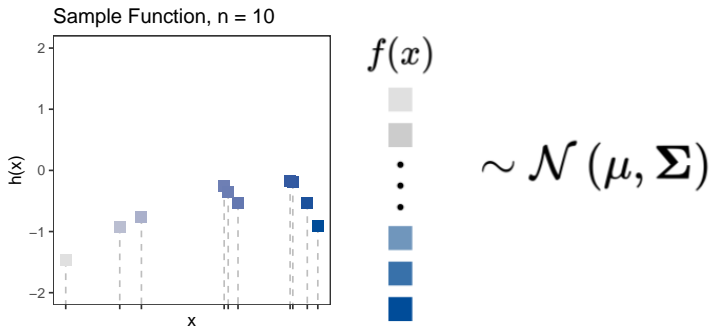$\sim \mathcal{N}\left(\mu, \boldsymbol{\Sigma}\right)$

GAUSSIAN PROCESSES: INTUITION

- Thus, it is required that for **any finite set** of inputs $\{\mathbf{x}^{(1)}, \ldots, \mathbf{x}^{(n)}\} \subset \mathcal{X}$, the vector **f** has a Gaussian distribution

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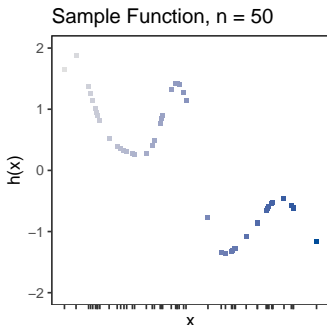Sample Function, n = 50

GAUSSIAN PROCESSES

This intuitive explanation is formally defined as follows:

A function $f(\mathbf{x})$ is generated by a GP $\mathcal{GP}\left(m(\mathbf{x}), k\left(\mathbf{x}, \mathbf{x}'\right)\right)$ if for **any finite** set of inputs $\left\{\mathbf{x}^{(1)}, \ldots, \mathbf{x}^{(n)}\right\}$, the associated vector of function values $\boldsymbol{f} = \left(f(\mathbf{x}^{(1)}), \ldots, f(\mathbf{x}^{(n)})\right)$ has a Gaussian distribution

$$\boldsymbol{f} = \left[f\left(\mathbf{x}^{(1)}\right), \ldots, f\left(\mathbf{x}^{(n)}\right)\right] \sim \mathcal{N}\left(\boldsymbol{m}, \boldsymbol{K}\right),$$

with

$$\mathbf{m} \ := \ \left(m\left(\mathbf{x}^{(i)}\right)\right)_i, \quad \mathbf{K} := \left(k\left(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}\right)\right)_{i,j},$$

where $m(\mathbf{x})$ is called mean function and $k(\mathbf{x}, \mathbf{x}')$ is called covariance function.

GAUSSIAN PROCESSES

A GP is thus **completely specified** by its mean and covariance function

$$
\begin{aligned}
m(\mathbf{x}) &= \mathbb{E}[f(\mathbf{x})] \\
k(\mathbf{x}, \mathbf{x}') &= \mathbb{E}\left[\left(f(\mathbf{x}) - \mathbb{E}[f(\mathbf{x})] \right) \left(f(\mathbf{x}') - \mathbb{E}[f(\mathbf{x}')] \right) \right]
\end{aligned}
$$

Note: For now, we assume $m(\mathbf{x}) \equiv 0$. This is not necessarily a drastic limitation - thus it is common to consider GPs with a zero mean function.

SAMPLING FROM A GAUSSIAN PROCESS PRIOR

We can draw functions from a Gaussian process prior. Let us consider
$f(\mathbf{x}) \sim \mathcal{GP}\left(0, k(\mathbf{x}, \mathbf{x}')\right)$ with the squared exponential covariance
function [*]

$$k(\mathbf{x}, \mathbf{x}') = \exp\left(-\frac{1}{2\ell^2}\|\mathbf{x} - \mathbf{x}'\|^2\right), \ \ell = 1.$$

This specifies the Gaussian process completely.

[*] We will talk later about different choices of covariance functions.

SAMPLING FROM A GAUSSIAN PROCESS PRIOR

To visualize a sample function, we

- choose a high number n (equidistant) points $\left\{ \mathbf{x}^{(1)}, \dots, \mathbf{x}^{(n)} \right\}$
- compute the corresponding covariance matrix
  $\mathbf{K} = \left(k\left(\mathbf{x}^{(i)}, \mathbf{x}^{(j)} \right) \right)_{i,j}$ by plugging in all pairs $\mathbf{x}^{(i)}, \mathbf{x}^{(j)}$
- sample from a Gaussian $\boldsymbol{f} \sim \mathcal{N}(\mathbf{0}, \boldsymbol{K})$.

We draw 10 times from the Gaussian, to get 10 different samples.
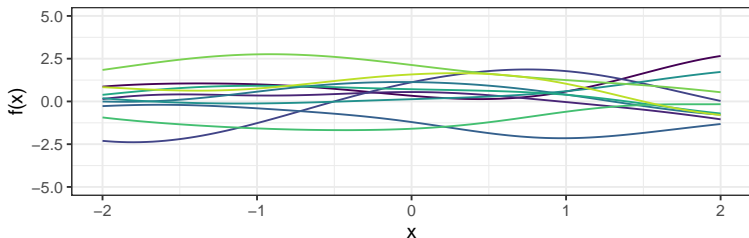
Since we specified the mean function to be zero $m(\mathbf{x}) \equiv 0$, the drawn functions have zero mean.

Gaussian Processes as Indexed Family

GAUSSIAN PROCESSES AS AN INDEXED FAMILY

A Gaussian process is a special case of a **stochastic process** which is defined as a collection of random variables indexed by some index set (also called an **indexed family**). What does it mean?

An **indexed family** is a mathematical function (or "rule") to map indices $t \in T$ to objects in $\mathcal{S}$.

Definition

A **family of elements in $\mathcal{S}$ indexed by** T (indexed family) is a surjective function

$$
\begin{aligned}
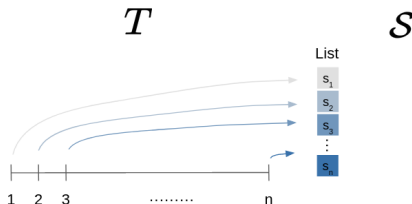s : T &\rightarrow \mathcal{S} \\
t &\mapsto s_t = s(t)
\end{aligned}
$$

INDEXED FAMILY

Some simple examples for indexed families are:
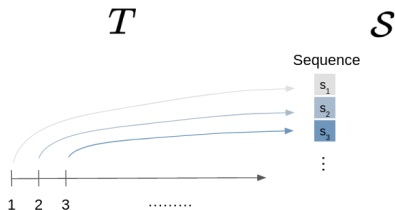
- finite sequences (lists): $T = \{1, 2, \ldots, n\}$ and $(s_t)_{t \in T} \in \mathbb{R}$
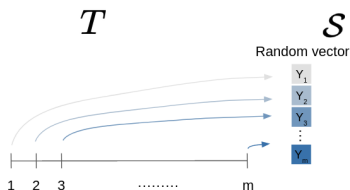

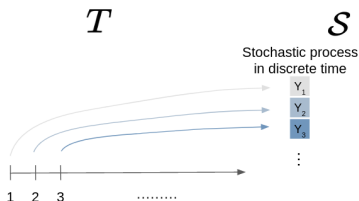- infinite sequences: $T = \mathbb{N}$ and $(s_t)_{t \in T} \in \mathbb{R}$

INDEXED FAMILY

But the indexed set $\mathcal{S}$ can be something more complicated, for example functions or **random variables** (RV):

- $T = \{1, \ldots, m\}$, Y_t's are RVs: Indexed family is a random vector.

- $T = \{1, \ldots, m\}$, Y_t's are RVs: Indexed family is a stochastic process in discrete time

- $T = \mathbb{Z}^2$, Y_t's are RVs: Indexed family is a 2D-random walk.

INDEXED FAMILY

- A Gaussian process is also an indexed family, where the random variables $f(\mathbf{x})$ are indexed by the input values $\mathbf{x} \in \mathcal{X}$.
- Their special feature: Any indexed (finite) random vector has a multivariate Gaussian distribution (which comes with all the nice properties of Gaussianity!).
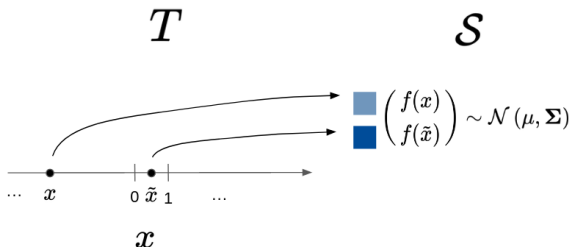
Visualization for a one-dimensional $\mathcal{X}$.

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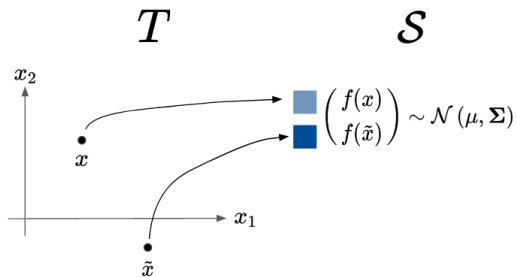
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